for net addition of claims) are hereby authorized to be charged to our Deposit Account No. 19-0036.

## Amendment

In the Claims:



Please cancel claims 11-14, 18-21, 26 and 27 without prejudice or disclaimer.

Please substitute the following claim 1 for the pending claim 1:

1. (Twice amended) A compound having the Formula I:

B2

or a pharmaceutically acceptable salt, prodrug or solvate thereof, wherein

X is one of O, S or CH<sub>2</sub>;

Het is

 $R_1$  is selected from the group consisting of hydrogen, optionally substituted alkyl, heteroaryl optionally substituted with one or more groups independently selected from the group consisting of halo, halo( $C_{1-6}$ )alkyl, hydroxy( $C_{1-6}$ )alkyl, amino( $C_{1-6}$ )alkyl, hydroxy, nitro,  $C_{1-6}$  alkyl,  $C_{1-6}$  alkoxy, aminocarbonyl, carbamoyloxy,  $C_{1-6}$  alkylsulfonylamino,  $C_{1-6}$  acyl and amino,  $C(O)R_{10}$ ,  $CH_2C(O)R_{10}$ ,  $S(O)R_{10}$ , and  $SO_2R_{10}$ ;

HOGENKAMP et al. Appl. No. 09/814,123

R<sub>2</sub> and R<sub>3</sub> are independently selected from the group consisting of hydrogen, alkyl, alkenyl, alkynyl, aryl, cyano, aminoalkyl, hydroxyalkyl, alkoxyalkyl, alkylthio, alkylsulfinyl, alkylsulfonyl, carboxyalkyl, alkylamino, dialkylamino, aminocarbonyl, alkylaminocarbonyl, arylaminocarbonyl, aralkylaminocarbonyl, alkylcarbonylamino, aralkylcarbonylamino, alkylcarbonyl, aminosulfonyl, alkylaminosulfonyl, and alkylsulfonyl;

R<sub>5</sub>, R<sub>6</sub>, R<sub>7</sub>, and R<sub>8</sub> are independently selected from the group consisting of hydrogen, halo, haloalkyl, alkyl, alkenyl, alkynyl, hydroxyalkyl, aminoalkyl, carboxyalkyl, alkoxyalkyl, nitro, amino, ureido, cyano, acylamino, amide, hydroxy, thiol, acyloxy, azido, alkoxy, carboxy, carbonylamido and alkylthiol;

R<sub>10</sub> is selected from the group consisting of amino, alkyl, alkenyl, alkynyl, OR<sub>11</sub>, alkylamino, dialkylamino, dialkylamino, dialkylaminoalkenyl, cycloalkyl, heterocycle, heteroaryl, aryl, aralkyl, arylalkenyl, arylalkynyl, and cycloalkylalkylamino, all of which are optionally substituted; and

 $R_{11}$  is selected from the group consisting of hydrogen, optionally substituted alkyl, and an alkalimetal.

Please substitute the following claim 15 for the pending claim 15:

15. (Twice Amended) A compound of claim 1, wherein:

 $R_1$  is  $C(O)R_{10}$ ,  $CH_2C(O)R_{10}$ , or  $SO_2R_{10}$ 

X is O or S;

 $R_{10}$  is amino, optionally substituted  $C_1$ - $C_6$  alkyl, or a heterocycle selected from the group consisting of N-morpholinyl, N-pyrrolidinyl and N-piperazinyl;

 $R_2$ , and  $R_3$  are independently hydrogen,  $C_1\text{-}C_6$  alkyl,  $C_1\text{-}C_6$  alkylthio or  $C_1\text{-}C_6$  alkylsulfinyl,

R<sub>5</sub> and R<sub>6</sub> are as defined in claim 1, and

 $R_7$  and  $R_8$  are independently selected from the group consisting of hydrogen, halo, halo( $C_1$ - $C_6$ )alkyl,  $C_1$ - $C_6$  alkyl, hydroxy( $C_1$ - $C_6$ )alkyl, amino( $C_1$ - $C_6$ )alkyl, carboxy( $C_1$ - $C_6$ )alkyl, aikoxy( $C_1$ - $C_6$ )alkyl, nitro, amino,  $C_1$ - $C_6$  acylamino, amide, hydroxy, thiol,  $C_1$ - $C_6$  acyloxy,  $C_1$ - $C_6$  alkoxy, carboxy, carbonylamido and  $C_1$ - $C_6$  alkylthiol.

Cont

B/3 Sub

I

16. (Twice Amended) A compound of Formula I:

or a pharmaceutically acceptable salt, prodrug/or solvate thereof, wherein

X is O or S;

Het is

$$-N$$
 $R_1$ 
 $R_2$ 

 $R_1$  is  $C(O)R_{10}$ ,  $CH_2C(O)R_{10}$ , or  $SO_2R_{10}$  wherein  $R_{10}$  is amino, alkyl, N-morpholinyl, N-pyrrolidinyl or N-piperazinyl, all of which are optionally substituted;

R<sub>2</sub> and R<sub>3</sub> are independently hydrogen, C<sub>1</sub>-C<sub>6</sub> alkyl, C<sub>1</sub>-C<sub>6</sub> alkylthio or C<sub>1</sub>-C<sub>6</sub> alkylsulfinyl; and

 $R_5$ ,  $R_6$ ,  $R_7$  and  $R_8$  are independently selected from the group consisting of hydrogen, halo, halo( $C_1$ - $C_6$ )alkyl,  $C_1$ - $C_6$  alkyl, hydroxy( $C_1$ - $C_6$ )alkyl, amino( $C_1$ - $C_6$ )alkyl, carboxy( $C_1$ - $C_6$ )alkyl, alkoxy( $C_1$ - $C_6$ )alkyl, nitro, amino,  $C_1$ - $C_6$  acylamino, amide, hydroxy, thiol,  $C_1$ - $C_6$  acyloxy,  $C_1$ - $C_6$  alkoxy, carboxy, carbonylamido and  $C_1$ - $C_6$  alkylthiol.

Sub (1)3

I

## 22. (Twice Amended) A compound of Formula 1:

R<sub>8</sub> R<sub>6</sub> Het

or a pharmaceutically acceptable salt, prodrug or solvate thereof, wherein

X is O or S;

Het is

-N R<sub>1</sub> R<sub>2</sub>

 $R_1$  is  $C(O)R_{10}$ , wherein  $R_{10}$  is amino, N-morpholinyl, N-pyrrolidinyl or N-piperazinyl, all of which are optionally substituted;

R<sub>2</sub> and R<sub>3</sub> are independently hydrogen, C<sub>1</sub>-C<sub>6</sub> alkyl, C<sub>1</sub>-C<sub>6</sub> alkylthio or C<sub>1</sub>-C<sub>6</sub> alkylsulfinyl; and

 $R_5$ ,  $R_6$ ,  $R_7$  and  $R_8$  are independently selected from the group consisting of hydrogen, halo, halo( $C_1$ - $C_6$ )alkyl,  $C_1$ - $C_6$  alkyl, hydroxy( $C_1$ - $C_6$ )alkyl, amino( $C_1$ - $C_6$ )alkyl, carboxy( $C_1$ - $C_6$ )alkyl, alkoxy( $C_1$ - $C_6$ )alkyl, nitro, amino,  $C_1$ - $C_6$  acylamino, amide, hydroxy, thiol,  $C_1$ - $C_6$  acyloxy,  $C_1$ - $C_6$  alkoxy, carboxy, carbonylamido and  $C_1$ - $C_6$  alkylthiol.

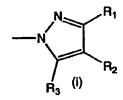
Sup.

I

## 25. (Once amended) A compound having the Formula I:

or a pharmaceutically acceptable salt, prodrug or solvate thereof, wherein

X is NR<sub>9</sub>C(O) or C(O)NR<sub>9</sub>, where R<sub>9</sub> is hydrogen or C<sub>1</sub>-C<sub>10</sub> alkyl; Het is



R<sub>1</sub> is SO<sub>2</sub>R<sub>10</sub>;

R<sub>2</sub> and R<sub>3</sub> are independently selected from the group consisting of hydrogen, alkyl, alkenyl, alkynyl, aryl, cyano, aminoalkyl, hydroxyalkyl, alkoxyalkyl, alkylthio, alkylsulfinyl, alkylsulfonyl, carboxyalkyl, alkylamino, dialkylamino, aminocarbonyl, alkylaminocarbonyl, arylaminocarbonyl, aralkylaminocarbonyl, alkylcarbonylamino, aralkylcarbonylamino, alkylcarbonyl, aminosulfonyl, alkylsulfonyl, and alkylsulfonyl;

R<sub>5</sub>, R<sub>6</sub>, R<sub>7</sub>, and R<sub>8</sub> are independently selected from the group consisting of hydrogen, halo, haloalkyl, alkyl, alkenyl, alkynyl, hydroxyalkyl, aminoalkyl, carboxyalkyl, alkoxyalkyl, nitro, amino, ureido, cyano, acylamino, amide, hydroxy, thiol, acyloxy, azido, alkoxy, carboxy, carbonylamido and alkylthiol;

R<sub>10</sub> is selected from the group consisting of amino, alkyl, alkenyl, alkynyl, OR<sub>11</sub>, alkylamino, dialkylamino, dialkylamino, dialkylaminoalkenyl, cycloalkyl, heterocycle, heteroaryl, aryl, aralkyl, arylalkenyl, arylalkynyl, and cycloalkylalkylamino, all of which are optionally substituted; and

R<sub>11</sub> is selected from the group consisting of hydrogen, optionally substituted alkyl, and an alkalimetal.

